

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF040219\
 Data File : BF113521.D
 Acq On : 2 Apr 2019 22:39
 Operator : JU/SJ
 Sample : K2180-19
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-10

Manual Integrations
 APPROVED

Sohil
 4/3/2019 12:42:49 PM

Quant Time: Apr 03 08:10:10 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	134862	20.00	ng	-0.01
21) Naphthalene-d8	7.97	136	472136	20.00	ng	-0.01
39) Acenaphthene-d10	9.72	164	221580	20.00	ng	-0.02
64) Phenanthrene-d10	11.20	188	363255	20.00	ng	-0.02
76) Chrysene-d12	13.83	240	332785	20.00	ng	-0.02
87) Perylene-d12	15.23	264	322066	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	872402	105.77	ng	0.00
7) Phenol-d6	6.35	99	1104133	105.82	ng	-0.01
23) Nitrobenzene-d5	7.26	82	652155	81.98	ng	-0.01
42) 2,4,6-Tribromophenol	10.51	330	275463	100.64	ng	-0.02
45) 2-Fluorobiphenyl	9.05	172	1193820	84.30	ng	-0.02
79) Terphenyl-d14	12.78	244	997591	66.07	ng	-0.02
Target Compounds						Qvalue
37) 2-Methylnaphthalene	8.69	142	40877	2.839	ng	97
38) 1-Methylnaphthalene	8.78	142	60726	4.385	ng	100
50) Dimethylphthalate	9.44	163	164502	10.018	ng	100
71) Phenanthrene	11.22	178	115940	6.427	ng	98
75) Fluoranthene	12.41	202	160946	7.892	ng	99
78) Pyrene	12.63	202	155690	6.751	ng	97
81) Benzo(a)anthracene	13.82	228	80777	4.240	ng	99
83) Chrysene	13.86	228	85177	4.402	ng	97
86) Indeno(1,2,3-cd)pyrene	16.59	276	54751	3.297	ng	98
88) Benzo(b)fluoranthene	14.84	252	137854m	6.890	ng	
89) Benzo(k)fluoranthene	14.86	252	40026m	2.278	ng	
90) Benzo(a)pyrene	15.18	252	86771	4.892	ng	# 96
92) Benzo(g,h,i)perylene	17.00	276	54116	3.500	ng	95

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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